

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081225\
 Data File : BF143349.D
 Acq On : 12 Aug 2025 12:07
 Operator : RC/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Quant Time: Aug 12 13:20:56 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 12 13:02:27 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.934	152	161554	20.000	ng	0.00
21) Naphthalene-d8	8.216	136	611608	20.000	ng	0.00
39) Acenaphthene-d10	9.969	164	320321	20.000	ng	0.00
64) Phenanthrene-d10	11.451	188	460173	20.000	ng	0.00
76) Chrysene-d12	14.092	240	273246	20.000	ng	0.00
86) Perylene-d12	15.586	264	337675	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	1115976	119.453	ng	0.00
7) Phenol-d6	6.569	99	1429946	119.357	ng	0.00
23) Nitrobenzene-d5	7.492	82	1251642	127.441	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	314370	134.290	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	2021003	106.990	ng	0.00
79) Terphenyl-d14	13.033	244	1568868	102.507	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.763	88	296653	61.364	ng	100
3) Pyridine	3.534	79	808240	62.740	ng	99
4) n-Nitrosodimethylamine	3.481	42	375923	62.645	ng	99
6) Aniline	6.592	93	996754	59.015	ng	# 88
8) 2-Chlorophenol	6.722	128	616100	61.717	ng	98
9) Benzaldehyde	6.481	77	387880	48.985	ng	100
10) Phenol	6.581	94	840504	59.221	ng	91
11) bis(2-Chloroethyl)ether	6.663	93	619725	59.247	ng	99
12) 1,3-Dichlorobenzene	6.875	146	671800	58.344	ng	99
13) 1,4-Dichlorobenzene	6.951	146	661936	57.278	ng	100
14) 1,2-Dichlorobenzene	7.104	146	634447	57.835	ng	100
15) Benzyl Alcohol	7.075	79	531507	61.891	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.198	45	1077772	57.714	ng	93
17) 2-Methylphenol	7.186	107	513134	60.895	ng	98
18) Hexachloroethane	7.445	117	225784	62.318	ng	100
19) n-Nitroso-di-n-propyla...	7.345	70	423943	57.269	ng	99
20) 3+4-Methylphenols	7.339	107	594674	58.354	ng	99
22) Acetophenone	7.339	105	765201	56.784	ng	98
24) Nitrobenzene	7.516	77	672780	63.075	ng	99
25) Isophorone	7.751	82	1195220	59.613	ng	99
26) 2-Nitrophenol	7.828	139	276412	63.401	ng	100
27) 2,4-Dimethylphenol	7.863	122	522594	62.476	ng	100
28) bis(2-Chloroethoxy)met...	7.957	93	726373	59.338	ng	99
29) 2,4-Dichlorophenol	8.075	162	512513	62.732	ng	99
30) 1,2,4-Trichlorobenzene	8.157	180	510824	59.275	ng	99
31) Naphthalene	8.239	128	1682766	57.179	ng	100
32) Benzoic acid	7.986	122	248714	62.282	ng	99
33) 4-Chloroaniline	8.281	127	632256	59.564	ng	99
34) Hexachlorobutadiene	8.357	225	314445	58.655	ng	99
35) Caprolactam	8.657	113	148587	64.696	ng	96
36) 4-Chloro-3-methylphenol	8.763	107	525032	61.301	ng	99
37) 2-Methylnaphthalene	8.928	142	1088275	56.249	ng	98
38) 1-Methylnaphthalene	9.028	142	1052909	56.505	ng	100
40) 1,2,4,5-Tetrachloroben...	9.092	216	516108	58.880	ng	99
41) Hexachlorocyclopentadiene	9.080	237	184941	61.485	ng	99
43) 2,4,6-Trichlorophenol	9.204	196	330076	66.369	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.245	196	355498	63.299	ng	98
46) 1,1'-Biphenyl	9.392	154	1282862	56.642	ng	99
47) 2-Chloronaphthalene	9.416	162	1038548	58.223	ng	99
48) 2-Nitroaniline	9.504	65	322909	61.917	ng	98
49) Acenaphthylene	9.828	152	1481001	57.494	ng	99
50) Dimethylphthalate	9.680	163	1125050	60.109	ng	99
51) 2,6-Dinitrotoluene	9.745	165	224724	62.496	ng	99
52) Acenaphthene	10.004	154	977067	57.055	ng	99
53) 3-Nitroaniline	9.916	138	270085	70.462	ng	99
54) 2,4-Dinitrophenol	10.022	184	82451	62.319	ng	# 77
55) Dibenzofuran	10.175	168	1389556	56.379	ng	99
56) 4-Nitrophenol	10.075	139	186307	62.130	ng	99
57) 2,4-Dinitrotoluene	10.151	165	295520	61.944	ng	95
58) Fluorene	10.516	166	1046138	55.570	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.292	232	265201	70.050	ng	99
60) Diethylphthalate	10.386	149	990716	59.094	ng	99
61) 4-Chlorophenyl-phenyle...	10.504	204	515436	56.664	ng	99
62) 4-Nitroaniline	10.527	138	242938	65.508	ng	98
63) Azobenzene	10.669	77	1104323	57.669	ng	98
65) 4,6-Dinitro-2-methylph...	10.563	198	119734	61.844	ng	98
66) n-Nitrosodiphenylamine	10.622	169	898967	59.328	ng	99
67) 4-Bromophenyl-phenylether	10.998	248	318696	60.555	ng	98
68) Hexachlorobenzene	11.069	284	341976	60.282	ng	100
69) Atrazine	11.145	200	250602	63.410	ng	98
70) Pentachlorophenol	11.263	266	165675	61.169	ng	98
71) Phenanthrene	11.480	178	1426848	56.831	ng	99
72) Anthracene	11.533	178	1459751	57.330	ng	100
73) Carbazole	11.680	167	1232939	56.857	ng	99
74) Di-n-butylphthalate	12.010	149	1127365	61.437	ng	100
75) Fluoranthene	12.669	202	1181812	53.737	ng	99
77) Benzidine	12.780	184	419198	56.115	ng	99
78) Pyrene	12.898	202	1196849	52.212	ng	100
80) Butylbenzylphthalate	13.504	149	283998	61.669	ng	98
81) Benzo(a)anthracene	14.080	228	1072649	61.763	ng	99
82) 3,3'-Dichlorobenzidine	14.039	252	313084	66.762	ng	99
83) Chrysene	14.121	228	937047	57.863	ng	100
84) Bis(2-ethylhexyl)phtha...	14.068	149	426553	62.688	ng	98
85) Di-n-octyl phthalate	14.680	149	856960	62.516	ng	99
87) Indeno(1,2,3-cd)pyrene	17.115	276	1462168	60.440	ng	100
88) Benzo(b)fluoranthene	15.145	252	1161807	62.905	ng	99
89) Benzo(k)fluoranthene	15.180	252	1035022	57.673	ng	99
90) Benzo(a)pyrene	15.521	252	1052580	60.510	ng	100
91) Dibenzo(a,h)anthracene	17.127	278	1170363	60.556	ng	99
92) Benzo(g,h,i)perylene	17.574	276	1182985	60.761	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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